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ULTRAVIOLET ABSORPTION CROSS-SECTIONS OF SOME CARBONYL COMPOUNDS AND THEIR TEMPERATURE DEPENDENCE.

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ABSTRACT

Ultraviolet absorption cross-section of phosgene (CCl_2O), trichloroacetylchloride ($\text{CCl}_3\text{-CClO}$) and trichloroacetaldehyde ($\text{CCl}_3\text{-CHO}$) have been measured between 170 and 320 nm for temperature ranging from 210 to 295 K with a classical double beam equipment.

These data are compared with other available determinations performed at room temperature. Photodissociation coefficients are estimated and their temperature dependence is discussed. Impact of the photodissociation on the total atmospheric destruction of these compounds is illustrated.

I. INTRODUCTION.

Chemical degradation of alternative hydrochlorofluorocarbons in troposphere produces a series of carbonyl compounds like phosgene, halo-aldehydes or halo-ketones, which could themselves be removed in three ways : (1) Photodissociation by solar UV radiations to produce potential odd chlorine precursors, (2) Reaction with H_2O and (3) Reaction with OH.

In order to determine the lifetimes for the photolysis processes, an accurate knowledge of the UV absorption cross-sections is required as a function of wavelength and temperature.

II. EXPERIMENTAL.

Ultraviolet absorption cross sections of phosgene (CCl_2O), trichloroacetylchloride ($\text{CCl}_3\text{-CClO}$) and trichloroacetaldehyde ($\text{CCl}_3\text{-CHO}$) have been measured between 170 and 320 nm for temperature ranging from 210 to 295 K with a classical double beam equipment. (Gillotay et al., 1989). The purity of the three compounds is better than 99.5 % as determined by gas phase chromatography.

III. RESULTS.

Numerical values of absorption cross-sections for wavenumber intervals of 500 cm^{-1} are given in tables I-III. The absorption spectra are illustrated in Figures 1-3 for selected temperatures namely 295, 250 and 210 K and compared with other available data at 295 K.

In all cases, Beer-Lambert's law was verified for absorption ranging from 10 to 85 %. In such conditions, and according to the error budget previously published, (Simon et al., 1988), the absorption cross-sections reported here are determined with an accuracy of $\pm 2 \%$ at room temperature and of ± 3 to $\pm 4 \%$ at the lowest temperature.

Carbonyl compounds display a continuous absorption in the 170-335 nm range. The presence of two maxima and the temperature dependence observed near the maxima and for the longest wavelengths, seems to indicate that there are two continua, one corresponding to the absorption of the C-Cl bond, with a maximum around 170 nm and the other centred respectively at 240 nm for phosgene, around 260 nm for trichloroacetylchloride and at 290 nm for trichloroacetaldehyde, corresponding to the absorption of the C=O bond.

Absorption cross-sections values change with temperature by a factor, which depends on both the wavelength and the chemical composition of the compound. For each wavelength, an exponential dependence of the absorption cross-section versus temperature is clearly established, with a decrease of absorption cross-sections in the region of low absorptions (up to 80 % at 305 nm and 210 K in the case of CCl_2O) and a small increase near the maximum of absorption (up to 5 %). This effect is the most important at the lowest temperature.

Discrepancies observed between the different set of available data at room temperature have to be discussed in more details in terms of experimental conditions.

IV. DISCUSSION.

Photodissociation coefficients of the molecules have been calculated, neglecting the effects of multiple scattering, for given altitude (z), zenith angle (χ) and wavelengths intervals according to the relation :

$$J(z) = \sigma_{\lambda} q_{\lambda}(z) \quad ; \quad q_{\lambda}(z) = q_{\lambda}(w) e^{-\tau_{\lambda}(z)}$$

$$\tau_{\lambda}(z) = \int_z^{\infty} [n(\text{O}_2) \sigma_{\lambda}(\text{O}_2) + n(\text{O}_3) \sigma_{\lambda}(\text{O}_3) + n(\text{air}) \sigma_{\text{scatt}}] \sec \chi \, dz$$

where

z is the altitude,

σ are the absorption cross-sections,

$q(z)$ and $q(\infty)$ are the solar irradiance at altitude z or extraterrestrial ($z = \infty$)

n is the number of particles per volume unit.

Calculations are made for solar zenith angle of 0° and 60° ($\sec = 1$ and 2), taking into account the values of $\sigma(O_2)$ and $\sigma(O_3)$ from WMO and Kockarts (1976), σ_{scatt} from Nicolet (1984) and the values of $q(\infty)$ from WMO (1986) and by taking into account the actual values of cross-sections corresponding to the temperature conditions at each altitude.

These photodissociation coefficients are illustrated in Figures 4-6 and compared with those calculated with values of absorption cross-sections measured at room temperature.

Stratospheric photodissociation coefficients (for altitude ranging from 15 to 50 km) calculated the temperature dependent absorption cross-sections, are smaller than those calculated with the room temperature values in the 20-35 km region, due to the decrease in the absorption cross-sections in the 200 nm region and the influence of the wavelengths longer than 280nm in the low stratosphere.

Tropospheric photodissociation coefficients for phosgene are very low (between 10^{-9} and $10^{-11} \text{ sec}^{-1}$) and are reduced down to 20 % of their room temperature values, using the temperature dependent cross-sections. For the two other compounds, tropospheric photodissociation coefficients are relatively high (between 10^{-7} and 10^{-4} sec^{-1}) and show a small temperature dependence. Photolysis is, for these two molecules, a non negligible mechanism for their tropospheric removal.

In conclusion, this work presents a new set of experimental data on the absorption cross-sections of carbonyl compounds in atmospheric temperature condition and highlights a non negligible temperature dependence of their photolysis.

V. REFERENCE.

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Table I PHOSGENE (CCl_2O)

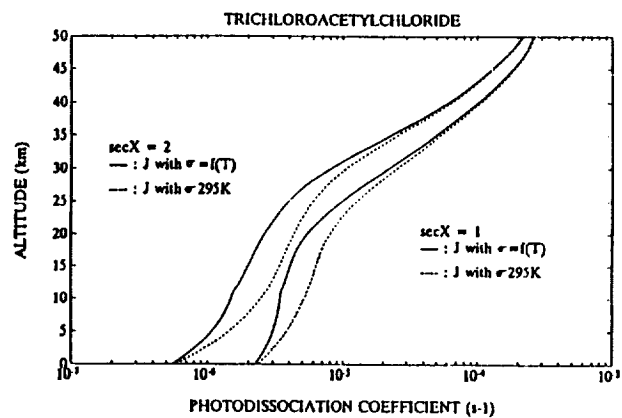
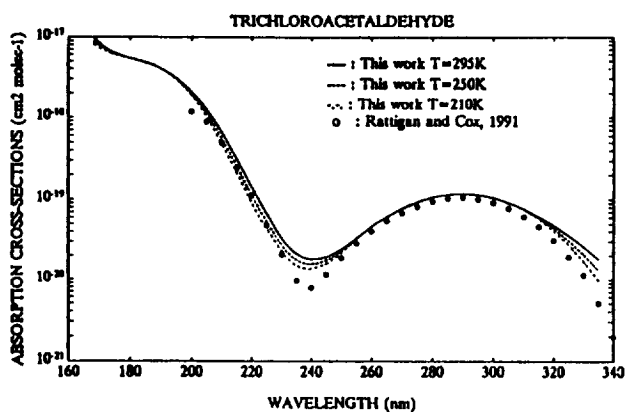
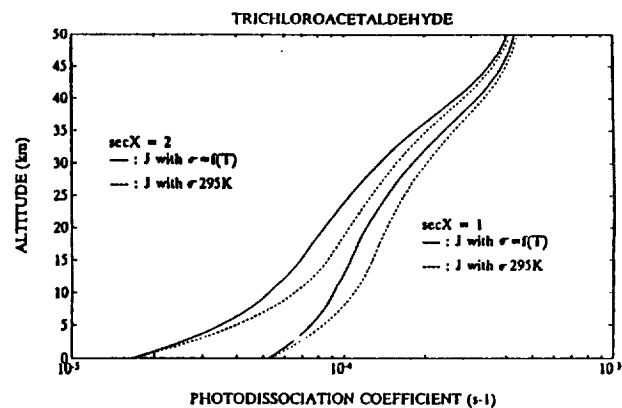
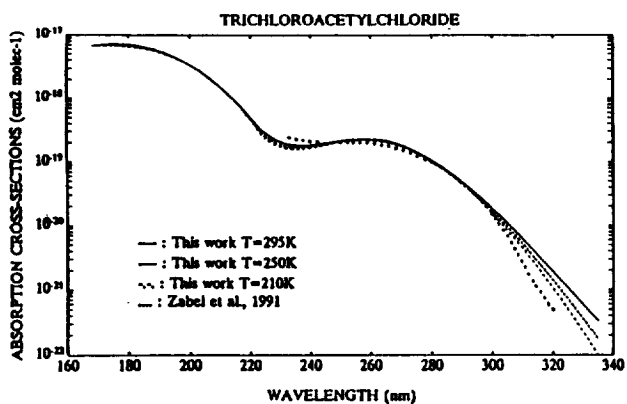
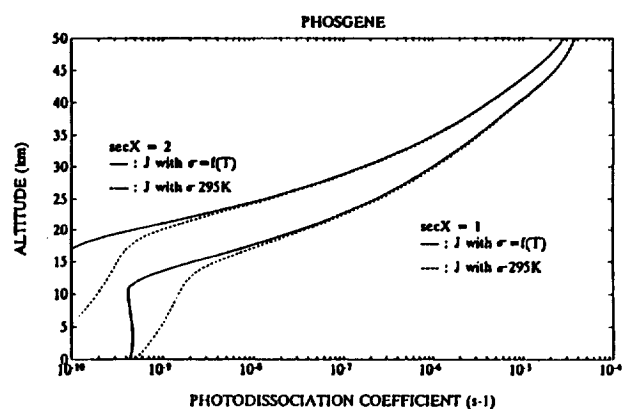
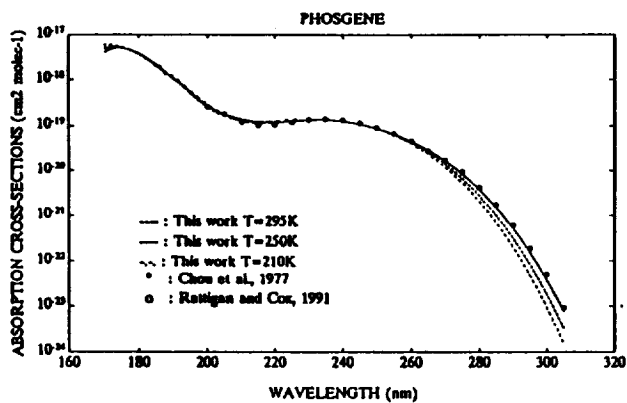
		$\sigma(\lambda) \times 10^{21} \text{ (cm}^2 \text{ molec}^{-1}\text{)}$				
λ	(nm)	295K	270K	250K	230K	210K
42	166.7-169.5	3010	2600	4140	4610	5560
43	169.5-172.4	4330	4760	5120	5530	5960
44	172.4-173.9	4930	5160	5360	5560	5770
45	173.2-175.4	5030	5190	5240	5330	5430
46	175.4-177.0	4890	4910	4920	4940	4950
47	177.0-178.6	4540	4490	4460	4420	4380
48	178.6-180.2	4040	3960	3900	3850	3790
49	180.2-181.8	3470	3390	3320	3270	3210
50	181.8-183.5	2890	2820	2760	2710	2660
51	183.5-185.2	2340	2280	2240	2200	2150
52	185.2-186.9	1860	1820	1790	1760	1730
53	186.9-188.7	1460	1440	1410	1380	1350
54	188.7-190.5	1160	1140	1120	1100	1080
55	190.5-192.3	903	888	876	865	854
56	192.3-194.2	715	698	685	673	660
57	194.2-196.1	524	509	498	487	476
58	196.1-198.0	399	384	376	367	357
59	198.0-200.0	312	301	293	285	277
60	200.0-202.0	252	243	236	229	223
61	202.0-204.1	209	202	196	191	185
62	204.1-206.2	179	173	168	163	159
63	206.2-208.3	158	153	149	145	141
64	208.3-210.5	143	139	136	132	129
65	210.5-212.8	123	120	117	114	111
66	212.8-215.0	106	104	101	98	95
67	215.0-217.4	92	90	87	84	81
68	217.4-219.6	81	79	76	73	70
69	219.6-221.2	72	70	67	64	61
70	221.2-223.7	64	62	59	56	53
71	223.7-226.9	57	55	52	49	46
72	226.9-229.9	51	49	46	43	40
73	229.9-232.6	46	44	41	38	35
74	232.6-235.3	41	39	36	33	30
75	235.3-238.1	37	35	32	29	26
76	238.1-241.0	33	31	28	25	22
77	241.0-243.9	30	28	25	22	19
78	243.9-246.9	27	25	22	19	16
79	246.9-250.0	24	22	19	16	13
80	250.0-253.2	21	19	16	13	10
81	253.2-256.4	19	17	14	11	8
82	256.4-259.7	17	15	12	9	6
83	259.7-263.2	15	13	10	7	4
84	263.2-266.7	13	11	8	5	3
85	266.7-270.3	11	9	6	4	2
86	270.3-274.0	10	8	5	3	1
87	274.0-277.8	9	7	4	2	1
88	277.8-281.7	8	6	3	1	0
89	281.7-285.7	7	5	2	1	0
90	285.7-289.9	6	4	1	0	0
91	289.9-294.1	5	3	1	0	0
92	294.1-298.5	4	2	0	0	0
93	298.5-303.0	3	1	0	0	0
94	303.0-307.7	2	1	0	0	0

Table III TRICHLOROACETYLCHLORIDE (CCl₃CClO)

N°	(nm)	$\epsilon(\lambda) \times 10^{21} \text{ (cm}^2 \text{ molec.}^{-1}\text{)}$				N°	(nm)	$\epsilon(\lambda) \times 10^{21} \text{ (cm}^2 \text{ molec.}^{-1}\text{)}$			
		295K	270K	250K	210K			295K	270K	250K	210K
42	166.7-169.5	6710	6710	6700	6690	42	166.7-169.5	9880	9210	8880	8070
43	169.5-172.4	6990	6910	6840	6720	43	169.5-172.4	7850	7540	7170	6470
44	172.4-175.3	7160	7000	6900	6810	44	172.4-175.3	6890	6710	6370	5660
45	175.3-178.2	7160	7020	6910	6800	45	175.3-178.2	6420	6300	6210	6120
46	178.2-181.1	7170	7010	6890	6770	46	178.2-181.1	6040	5980	5920	5870
47	181.1-184.0	7140	6970	6840	6710	47	181.1-184.0	5720	5700	5680	5630
48	184.0-186.9	7060	6890	6750	6620	48	184.0-186.9	5480	5480	5470	5470
49	186.9-189.8	6940	6770	6630	6500	49	186.9-189.8	5250	5270	5290	5320
50	189.8-192.7	6770	6600	6460	6330	50	189.8-192.7	5030	5070	5090	5150
51	192.7-195.6	6550	6380	6240	6110	51	192.7-195.6	4820	4850	4880	4930
52	195.6-198.5	6320	6150	6010	5880	52	195.6-198.5	4620	4620	4650	4710
53	198.5-201.4	6100	5930	5790	5660	53	198.5-201.4	4390	4390	4410	4430
54	201.4-204.3	5880	5710	5570	5440	54	201.4-204.3	4100	4080	4100	4120
55	204.3-207.2	5660	5490	5350	5220	55	204.3-207.2	3740	3740	3740	3740
56	207.2-210.1	5440	5270	5130	5000	56	207.2-210.1	3360	3360	3360	3360
57	210.1-213.0	5220	5050	4910	4780	57	210.1-213.0	2990	2990	2990	2990
58	213.0-215.9	5000	4830	4690	4560	58	213.0-215.9	2650	2650	2650	2650
59	215.9-218.8	4780	4610	4470	4340	59	215.9-218.8	2200	2200	2200	2200
60	218.8-221.7	4560	4390	4250	4120	60	218.8-221.7	1750	1750	1750	1750
61	221.7-224.6	4340	4170	4030	3900	61	221.7-224.6	1390	1390	1390	1390
62	224.6-227.5	4120	3950	3810	3680	62	224.6-227.5	1080	1080	1080	1080
63	227.5-230.4	3900	3730	3590	3460	63	227.5-230.4	812	812	812	812
64	230.4-233.3	3680	3510	3370	3240	64	230.4-233.3	596	596	596	596
65	233.3-236.2	3460	3290	3150	3020	65	233.3-236.2	424	424	424	424
66	236.2-239.1	3240	3070	2930	2800	66	236.2-239.1	298	298	298	298
67	239.1-242.0	3020	2850	2710	2580	67	239.1-242.0	203	203	203	203
68	242.0-244.9	2800	2630	2490	2360	68	242.0-244.9	139	139	139	139
69	244.9-247.8	2580	2410	2270	2140	69	244.9-247.8	85.7	85.7	85.7	85.7
70	247.8-250.7	2360	2190	2050	1920	70	247.8-250.7	51.2	51.2	51.2	51.2
71	250.7-253.6	2140	1970	1830	1700	71	250.7-253.6	35.4	35.4	35.4	35.4
72	253.6-256.5	1920	1750	1610	1480	72	253.6-256.5	24.8	24.8	24.8	24.8
73	256.5-259.4	1700	1530	1390	1260	73	256.5-259.4	17.5	17.5	17.5	17.5
74	259.4-262.3	1480	1310	1170	1040	74	259.4-262.3	12.2	12.2	12.2	12.2
75	262.3-265.2	1260	1090	950	820	75	262.3-265.2	8.4	8.4	8.4	8.4
76	265.2-268.1	1040	870	730	600	76	265.2-268.1	5.1	5.1	5.1	5.1
77	268.1-271.0	820	650	510	380	77	268.1-271.0	3.3	3.3	3.3	3.3
78	271.0-273.9	600	430	290	160	78	271.0-273.9	2.1	2.1	2.1	2.1
79	273.9-276.8	380	210	70	30	79	273.9-276.8	1.4	1.4	1.4	1.4
80	276.8-279.7	160	90	30	10	80	276.8-279.7	0.9	0.9	0.9	0.9
81	279.7-282.6	90	50	20	10	81	279.7-282.6	0.6	0.6	0.6	0.6
82	282.6-285.5	50	30	10	5	82	282.6-285.5	0.4	0.4	0.4	0.4
83	285.5-288.4	30	20	10	5	83	285.5-288.4	0.3	0.3	0.3	0.3
84	288.4-291.3	20	10	5	5	84	288.4-291.3	0.2	0.2	0.2	0.2
85	291.3-294.2	10	5	5	5	85	291.3-294.2	0.1	0.1	0.1	0.1
86	294.2-297.1	5	5	5	5	86	294.2-297.1	0.1	0.1	0.1	0.1
87	297.1-300.0	5	5	5	5	87	297.1-300.0	0.1	0.1	0.1	0.1
88	300.0-302.9	5	5	5	5	88	300.0-302.9	0.1	0.1	0.1	0.1
89	302.9-305.8	5	5	5	5	89	302.9-305.8	0.1	0.1	0.1	0.1
90	305.8-308.7	5	5	5	5	90	305.8-308.7	0.1	0.1	0.1	0.1
91	308.7-311.6	5	5	5	5	91	308.7-311.6	0.1	0.1	0.1	0.1
92	311.6-314.5	5	5	5	5	92	311.6-314.5	0.1	0.1	0.1	0.1
93	314.5-317.4	5	5	5	5	93	314.5-317.4	0.1	0.1	0.1	0.1
94	317.4-320.3	5	5	5	5	94	317.4-320.3	0.1	0.1	0.1	0.1
95	320.3-323.2	5	5	5	5	95	320.3-323.2	0.1	0.1	0.1	0.1
96	323.2-326.1	5	5	5	5	96	323.2-326.1	0.1	0.1	0.1	0.1
97	326.1-329.0	5	5	5	5	97	326.1-329.0	0.1	0.1	0.1	0.1
98	329.0-331.9	5	5	5	5	98	329.0-331.9	0.1	0.1	0.1	0.1
99	331.9-334.8	5	5	5	5	99	331.9-334.8	0.1	0.1	0.1	0.1
100	334.8-337.7	5	5	5	5	100	334.8-337.7	0.1	0.1	0.1	0.1

Table II TRICHLOROACETALDEHYDE (CCl₃CHO)

$\epsilon(\lambda) \times 10^{21} \text{ (cm}^2 \text{ molec.}^{-1}\text{)}$					
295K	270K	250K	230K	210K	
8880	9210	8880	8470	8070	
7550	7440	7210	7080	6860	
6990	6700	6710	6430	6300	
6200	6200	6100	5920	5820	
5980	5940	5930	5720	5620	
5720	5700	5680	5470	5370	
5480	5480	5470	5290	5200	
5250	5270	5290	5130	5070	
5030	5070	5090	4920	4930	
4820	4850	4880	4680	4710	
4620	4620	4650	4480	4470	
4390	4390	4410	4110	4120	
4100	4080	4100	3740	3740	
3740	3740	3740	3360	3330	
3360	3360	3360	2960	2900	
2990	2990	2990	2550	2460	
2650	2650	2650	2140	2090	
2200	2200	2200	1790	1760	
1750	1750	1750	1390	1380	
1390	1390	1390	1080	1020	
1080	1080	1080	812	762	
865	865	812	762	715	
703	642	596	554	515	
603	461	424	390	359	
366	326	298	271	248	
257	227	205	186	168	
176	155	139	126	113	
121	106	95.3	85.7	77.0	
82.5	73.2	65.9	59.4	53.5	
57.9	51.3	46.5	42.1	38.2	
38.4	35.3	32.1	29.4	26.8	
28.3	25.3	23.2	21.2	19.4	
22.3	20.1	18.4	17.0	15.6	
15.2	17.4	16.1	14.9	13.8	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
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16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
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18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
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18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
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18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
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16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4	18.4	15.4	14.3	13.4	
18.4	16.4	15.4	14.3	13.4	
16.4</					



Figures 1-3 : Ultraviolet absorption cross-sections at 295 K, 250 K and 210 K.

Figures 4-6 : Photodissociation coefficients as a function of altitude.